WHAT IS CLAIMED IS:

1. An indene derivative of formula (I) or a pharmaceutically acceptable salt thereof:

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(I)

wherein,

R_{1a} is OH or H;

R_{1b} is C₁₋₆ alkyl, C₃₋₆ cycloalkyl, benzyl or phenyl, the phenyl being optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂ and OR^a, when R_{1a} is OH; when R_{1a} is H,

R₂ is CN, CO₂R^a or CONR^eR^f;

 R_3 is phenyl optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a and C₁₋₆ alkyl; and

 R^4 , R^5 , R^6 and R^7 are each independently H, $O(CH_2)_m R^g$ or $CH_2 R^h$; in which

 R^a is H, C_{1-6} alkyl or C_{3-6} cycloalkyl, the C_{1-6} alkyl and C_{3-6} cycloalkyl being optionally substituted with one or more halogens;

 R^b , R^c , R^e and R^f are each independently H, C_{1-6} alkyl, C_{3-6} cycloalkyl or benzyl;

R^d is O, S or NR^a;

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R^g is H, , or phenyl, the phenyl being optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂ and NO₂;

$$R_h\,is$$
 -{-N-Rd}; and

m is an integer in the range of 1 to 3.

2. The compound of claim 1, wherein R_{1b} is C_{1-6} alkyl, C_{3-6} cycloalkyl, benzyl or phenyl, the phenyl being optionally substituted with one or more methoxy groups,

when R_{1a} is OH; when R_{1a} is H, R_{1b} is OR^a, NR^bR^c, NHCOR^a or ; R_3 is phenyl being optionally substituted with one or more halogens or C_{1-4} alkyls; and R_4 and R_7 is H, in which R^a is H or C_{1-6} alkyl; R^d is O or S; R^g is H, phenyl,

$$-\frac{1}{2} \frac{1}{|\nabla|_{N}} - \frac{1}{2} \frac{1}{|\nabla|_{N}}$$
 or

- 3. The compound of claim 1, wherein R_3 is phenyl, R_5 is H, and R_6 is $O(CH_2)_m R^g$ or CH_2R^h .
 - 4. The compound of claim 1, which is selected from the group consisting of:

 1-hydroxy-6-methoxy-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-methoxy-1-(3-methoxy-phenyl)-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1-isopropyl-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-methoxy-1-methyl-3-phenyl-1H-indene-2-carboxylic acid

ethyl ester,

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1-benzyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-cyclohexyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-(2-morpholine-4-yl-ethoxy)-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-morpholine-4-yl-methyl-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(2-pyridine-2-yl-ethoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-Carbonitrile,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid methyl ester,

1-hydroxy-6-methoxy-1,3-diphenyl-1H-indene-2-carboxylic acid,

1-hydroxy-6-methoxy-1-methyl-3-phenyl-1H-indene-2-carboxylic acid,

1-benzyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid, 1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-carboxylic acid,

1-cyclohexyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid,

1,6-dimethoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-ethoxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-amino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-amino-6-(2-morpholin-4-yl-ethoxy)-3-phenyl-1H-indene-2-carboxylic acid cyclohexyl amide,

1-amino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carbonitrile,

1-acetylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

6-methoxy-3-phenyl-1-propionylamino-1H-indene-2-carboxylic acid ethyl ester,

1-acetylamino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-acetylamino-6-(2-morpholin-4-yl-ethoxy)-3-phenyl-1H-indene-2carboxylic acid cyclohexyl amide,

1-diethylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester, 10

1-ethylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester, 6-methoxy-1-morpholin-4-yl-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-benzyl amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester, and 15

1-cyclohexyl amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester.

5. A process for preparing a compound of formula (I-a) which comprises reacting a compound of formula (II) with a Grignard reagent: 20

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wherein R_{1a} is OH; R_{1b} is alkyl, phenyl or benzyl; and R₂, R₃, R₄, R₅, R₆ and R₇ have

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the same meaning as defined in claim 1.

6. A process for preparing a compound of formula (I-e) which comprises reacting a compound of formula (II) with hydroxyl amine to obtain a compound of formula (III), and hydrogenation of the compound of formula (III) followed by reacting with acetyl chloride or an anhydrous acetic acid:

$$\begin{array}{c} R_{1b} \\ R_{2} \\ R_{3} \end{array} \qquad (I-e)$$

$$R_{5}$$
 R_{7} R_{8} R_{8} (II)

$$\begin{array}{c} HQ \\ HQ \\ R_{7} & N \\ R_{8} & R_{2} \end{array} \tag{IIII}$$

wherein R_{1a} is H; R_{1b} is NH₂ or NHCOR^a; and R₂, R₃, R₄, R₅, R₆ and R₇ have the same meaning as defined in claim 1.

7. A process for preparing a compound of formula (I-d) which comprises halogenation of a compound of formula (VIII) to obtain a compound of formula (IV), and reacting the compound of formula (IV) with an amine or alcohol compound:

wherein R_{1a} is H; R_{1b} is OR^a , NR^bR^c or R^d ; X is halogen; and R_2 , R_3 , R_4 , R_5 , R_6 and R_7 have the same meaning as defined in claim 1.

- 8. A pharmaceutical composition for modulating the activities of peroxisome proliferator activated receptors (PPARs) comprising a therapeutically effective amount of the compound or a salt defined in claim 1 as an active ingredient together with a pharmaceutically acceptable carrier.
- 9. The composition of claim 8, which is used for the treatment and prevention of diabetes, obesity, arteriosclerosis, hyperlipidemia, hyperinsulinism, hypertension, osteoporosis, liver cirrhosis, asthma and cancer.

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